



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 01:43 am BST

PDB ID : 6KIN
Title : Crystal structure of the tri-functional malyl-CoA lyase from *Roseiflexus castenholzii*
Authors : Tang, W.R.; Zhang, C.Y.; Wang, C.; Xu, X.L.
Deposited on : 2019-07-19
Resolution : 2.53 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

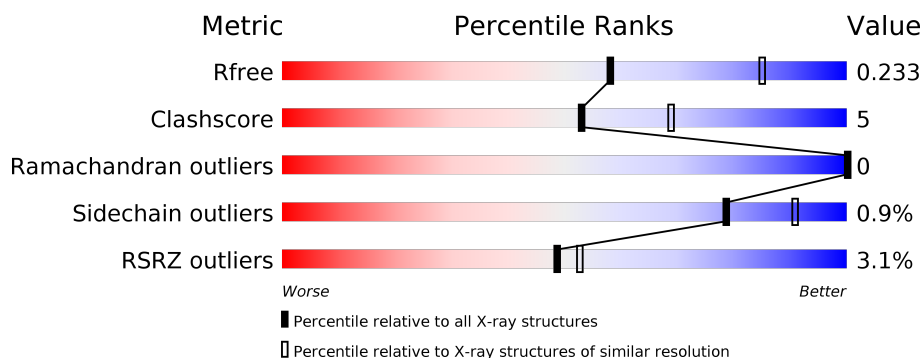
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.53 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5743 (2.54-2.50)
Clashscore	141614	6463 (2.54-2.50)
Ramachandran outliers	138981	6335 (2.54-2.50)
Sidechain outliers	138945	6337 (2.54-2.50)
RSRZ outliers	127900	5630 (2.54-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	347	<div> <div>2%</div> <div>88% 11% .</div> </div>
1	B	347	<div> <div>6%</div> <div>85% 14% ..</div> </div>
1	C	347	<div> <div>3%</div> <div>88% 11% .</div> </div>
1	D	347	<div> <div>%</div> <div>89% 10% .</div> </div>
1	E	347	<div> <div>3%</div> <div>83% 16%</div> </div>
1	F	347	<div> <div>2%</div> <div>85% 14% .</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	TRS	C	401	-	X	-	-

2 Entry composition [i](#)

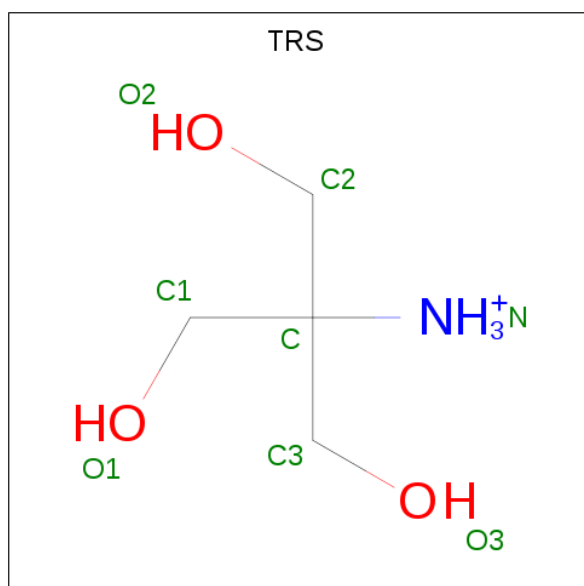
There are 3 unique types of molecules in this entry. The entry contains 16325 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called HpcH/HpaI aldolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	344	Total	C	N	O	S	0	0	0
			2670	1718	467	470	15			
1	B	344	Total	C	N	O	S	0	0	0
			2674	1720	467	472	15			
1	C	341	Total	C	N	O	S	0	0	0
			2646	1704	461	466	15			
1	D	345	Total	C	N	O	S	0	0	0
			2683	1725	469	474	15			
1	E	346	Total	C	N	O	S	0	0	0
			2689	1727	470	477	15			
1	F	342	Total	C	N	O	S	0	0	0
			2654	1709	461	469	15			

- Molecule 2 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			8	4	1	3		
2	C	1	Total	C	N	O	0	0
			8	4	1	3		

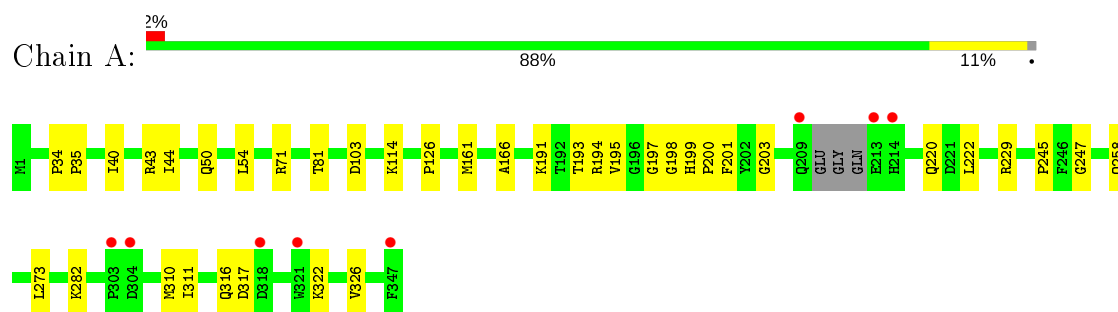
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	57	Total	O	0	0
			57	57		
3	B	49	Total	O	0	0
			49	49		
3	C	50	Total	O	0	0
			50	50		
3	D	59	Total	O	0	0
			59	59		
3	E	38	Total	O	0	0
			38	38		
3	F	40	Total	O	0	0
			40	40		

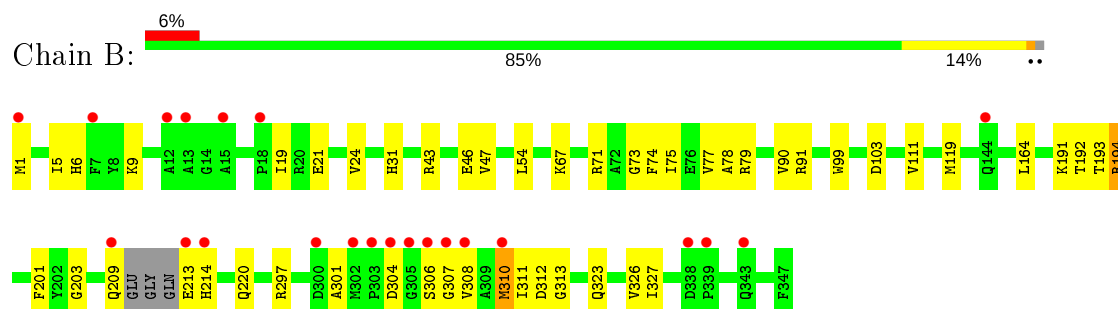
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

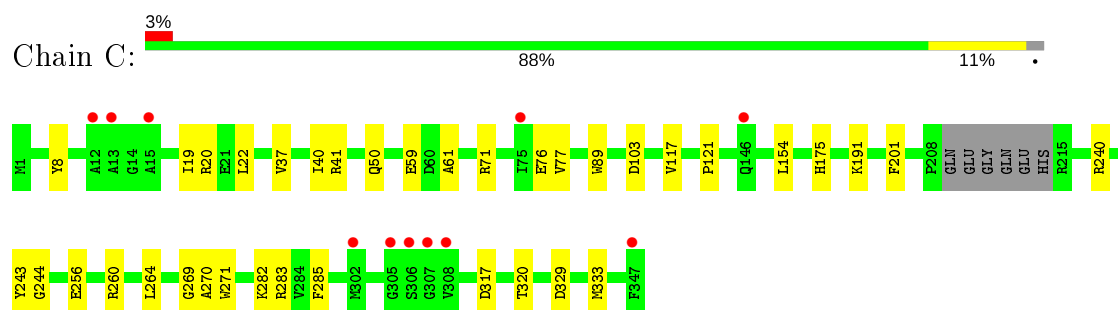
- Molecule 1: HpCh/HpaI aldolase



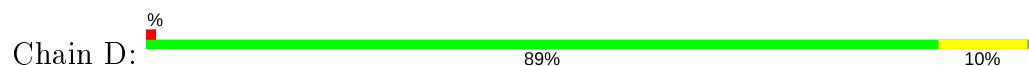
- Molecule 1: HpCh/HpaI aldolase

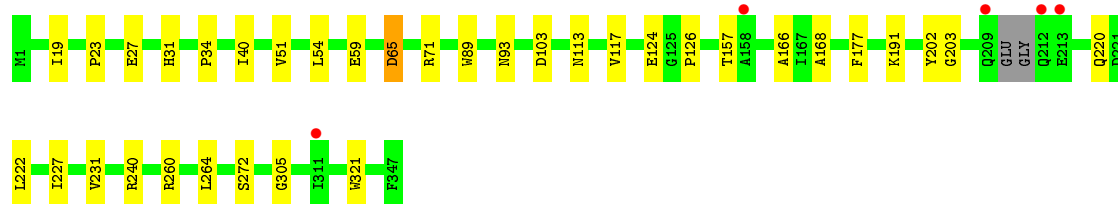


- Molecule 1: HpCh/HpaI aldolase

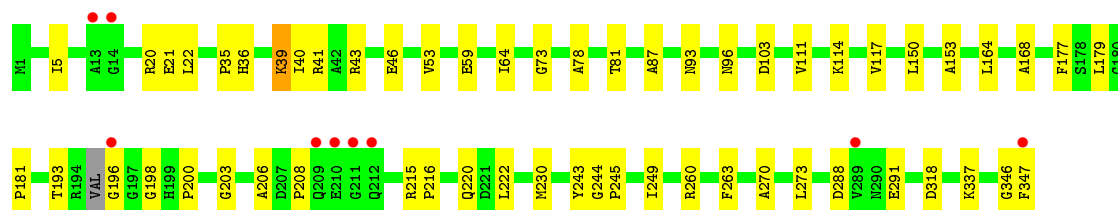
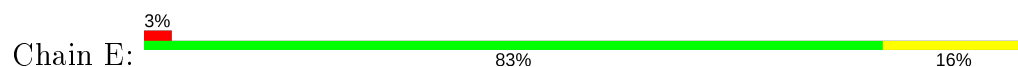


- Molecule 1: HpCh/HpaI aldolase

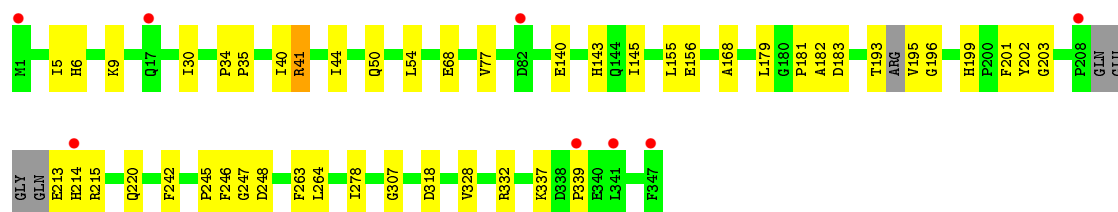
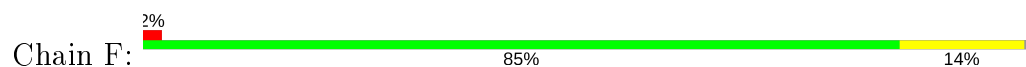




- Molecule 1: HpcH/HpaI aldolase



- Molecule 1: HpcH/HpaI aldolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	122.03Å 133.42Å 139.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.23 – 2.53 48.23 – 2.53	Depositor EDS
% Data completeness (in resolution range)	95.1 (48.23-2.53) 95.2 (48.23-2.53)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.176 , 0.234 0.177 , 0.233	Depositor DCC
R_{free} test set	3618 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	48.7	Xtriage
Anisotropy	0.290	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.017 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16325	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.10% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: TRS

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.46	0/2739	0.61	0/3722
1	B	0.44	0/2743	0.58	0/3727
1	C	0.47	0/2714	0.60	0/3688
1	D	0.43	0/2752	0.57	0/3739
1	E	0.50	0/2758	0.59	0/3746
1	F	0.44	0/2722	0.57	0/3698
All	All	0.46	0/16428	0.59	0/22320

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2670	0	2675	25	0
1	B	2674	0	2679	38	0
1	C	2646	0	2658	28	0
1	D	2683	0	2687	20	0
1	E	2689	0	2687	44	0
1	F	2654	0	2657	31	0
2	B	8	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	8	0	12	0	0
3	A	57	0	0	2	0
3	B	49	0	0	2	0
3	C	50	0	0	1	0
3	D	59	0	0	2	0
3	E	38	0	0	0	0
3	F	40	0	0	1	0
All	All	16325	0	16067	174	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (174) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:41:ARG:NH2	1:E:73:GLY:HA3	1.87	0.88
1:A:310:MET:HE3	1:E:64:ILE:HG13	1.56	0.85
1:C:41:ARG:HD2	1:C:77:VAL:HG23	1.60	0.83
1:B:74:PHE:CZ	1:B:90:VAL:HG22	2.15	0.81
1:A:310:MET:HE3	1:E:64:ILE:CG1	2.10	0.81
1:F:181:PRO:HB2	1:F:196:GLY:HA2	1.64	0.80
1:B:213:GLU:HG2	1:B:214:HIS:H	1.49	0.78
1:C:260:ARG:NH1	3:C:501:HOH:O	2.20	0.74
1:A:50:GLN:O	1:A:282:LYS:NZ	2.20	0.72
1:C:201:PHE:HB2	1:F:264:LEU:HD11	1.73	0.70
1:A:81:THR:O	1:A:114:LYS:NZ	2.24	0.70
1:E:41:ARG:NH2	1:E:73:GLY:CA	2.54	0.70
1:F:50:GLN:HG2	1:F:278:ILE:HG21	1.75	0.69
1:B:213:GLU:CG	1:B:214:HIS:H	2.05	0.69
1:B:193:THR:H	1:B:194:ARG:HH12	1.39	0.68
1:B:310:MET:HG3	1:B:310:MET:O	1.93	0.67
1:E:179:LEU:HB2	1:E:230:MET:HE1	1.77	0.65
1:B:192:THR:HB	1:B:194:ARG:HH22	1.61	0.63
1:E:249:ILE:HD11	1:E:273:LEU:HB2	1.80	0.63
1:E:288:ASP:HB2	1:E:291:GLU:HG3	1.82	0.62
1:A:193:THR:HG21	1:A:200:PRO:HD2	1.81	0.62
1:E:181:PRO:HB2	1:E:196:GLY:HA2	1.83	0.61
1:F:203:GLY:HA2	1:F:220:GLN:HG2	1.82	0.61
1:E:35:PRO:O	1:E:41:ARG:NE	2.30	0.60
1:E:193:THR:HG21	1:E:200:PRO:HD2	1.83	0.60
1:F:246:PHE:CE2	1:F:248:ASP:HB3	2.35	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:74:PHE:CZ	1:B:90:VAL:CG2	2.86	0.59
1:F:156:GLU:HG2	1:F:183:ASP:HB2	1.85	0.59
1:B:5:ILE:HB	3:B:511:HOH:O	2.03	0.59
1:D:65:ASP:OD1	1:D:65:ASP:N	2.37	0.58
1:D:51:VAL:HG21	1:D:54:LEU:HD13	1.83	0.58
1:F:199:HIS:CE1	1:F:201:PHE:HB2	2.38	0.58
1:A:317:ASP:OD2	1:E:96:ASN:ND2	2.35	0.58
1:B:191:LYS:HG2	1:E:263:PHE:CZ	2.40	0.57
1:B:193:THR:H	1:B:194:ARG:NH1	2.03	0.57
1:F:213:GLU:HG2	1:F:214:HIS:H	1.69	0.57
1:B:301:ALA:HB2	1:B:311:ILE:HD11	1.87	0.57
1:D:203:GLY:HA2	1:D:220:GLN:HG2	1.87	0.57
1:E:41:ARG:HH21	1:E:73:GLY:HA3	1.68	0.57
1:B:91:ARG:HA	1:B:119:MET:HB3	1.87	0.56
1:E:181:PRO:HB2	1:E:196:GLY:CA	2.36	0.56
1:F:5:ILE:HD12	1:F:6:HIS:CE1	2.41	0.56
1:A:191:LYS:HA	3:A:408:HOH:O	2.04	0.55
1:B:310:MET:HE1	1:B:313:GLY:HA2	1.88	0.55
1:B:73:GLY:O	1:B:77:VAL:HG23	2.05	0.55
1:A:193:THR:CG2	1:A:200:PRO:HD2	2.37	0.54
1:E:206:ALA:HB3	1:E:216:PRO:HD2	1.88	0.54
1:E:203:GLY:HA2	1:E:220:GLN:HG2	1.89	0.54
1:F:181:PRO:HD2	1:F:245:PRO:HD2	1.87	0.54
1:E:53:VAL:HG22	1:E:87:ALA:HB3	1.90	0.54
1:C:20:ARG:HH21	1:E:20:ARG:NH1	2.06	0.53
1:C:191:LYS:HG2	1:F:263:PHE:CZ	2.43	0.53
1:B:67:LYS:HG3	1:B:99:TRP:CE2	2.44	0.53
1:B:191:LYS:HG2	1:E:263:PHE:CE2	2.44	0.52
1:C:50:GLN:O	1:C:282:LYS:NZ	2.37	0.51
1:B:323:GLN:O	1:B:327:ILE:HG22	2.11	0.51
1:F:143:HIS:HB2	1:F:145:ILE:CD1	2.41	0.51
1:B:24:VAL:HG12	1:B:326:VAL:HG13	1.92	0.50
1:E:181:PRO:HD3	1:E:245:PRO:HD2	1.92	0.50
1:A:44:ILE:HG23	1:A:54:LEU:HD21	1.93	0.50
1:F:34:PRO:HD2	1:F:40:ILE:HG21	1.92	0.50
1:E:346:GLY:O	1:E:347:PHE:HB2	2.12	0.50
1:C:61:ALA:HA	1:F:307:GLY:O	2.11	0.50
1:F:30:ILE:HD11	1:F:242:PHE:CD2	2.47	0.50
1:A:195:VAL:HB	1:A:247:GLY:HA3	1.94	0.49
1:A:197:GLY:HA2	1:A:258:GLN:OE1	2.12	0.49
1:A:245:PRO:HB2	1:A:273:LEU:HD11	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:310:MET:HE1	1:B:313:GLY:CA	2.41	0.49
1:C:22:LEU:HD22	1:C:240:ARG:NH1	2.27	0.49
1:B:203:GLY:HA2	1:B:220:GLN:HG2	1.93	0.49
1:C:243:TYR:O	1:C:270:ALA:HA	2.13	0.49
1:E:168:ALA:HB2	1:E:177:PHE:HZ	1.77	0.49
1:D:168:ALA:HB2	1:D:177:PHE:HZ	1.77	0.49
1:A:161:MET:O	1:A:229:ARG:NH1	2.43	0.49
1:C:22:LEU:HD21	1:C:175:HIS:CD2	2.48	0.49
1:B:213:GLU:CG	1:B:214:HIS:N	2.73	0.48
1:E:208:PRO:HD3	1:E:215:ARG:NH2	2.28	0.48
1:E:93:ASN:ND2	1:E:103:ASP:OD2	2.31	0.48
1:D:113:ASN:ND2	3:D:401:HOH:O	2.30	0.48
1:F:195:VAL:HB	1:F:247:GLY:HA3	1.96	0.48
1:D:89:TRP:CD1	1:D:117:VAL:HB	2.49	0.47
1:D:34:PRO:HG2	1:D:40:ILE:HD13	1.96	0.47
1:C:329:ASP:O	1:C:333:MET:HG3	2.15	0.47
1:F:140:GLU:HA	1:F:145:ILE:HD13	1.96	0.47
1:F:199:HIS:HE1	1:F:201:PHE:HB2	1.78	0.47
1:C:71:ARG:HD2	1:C:103:ASP:HA	1.96	0.47
1:C:8:TYR:OH	1:E:5:ILE:HD13	2.15	0.47
1:F:6:HIS:HA	1:F:9:LYS:HD2	1.97	0.47
1:B:194:ARG:NH1	1:B:194:ARG:H	2.13	0.47
1:A:199:HIS:CE1	1:A:201:PHE:HB2	2.50	0.47
1:D:27:GLU:OE1	1:D:240:ARG:NH1	2.48	0.47
1:C:37:VAL:HG12	1:C:40:ILE:HG13	1.97	0.47
1:C:264:LEU:HD22	1:D:191:LYS:HB2	1.96	0.47
1:F:202:TYR:CE2	1:F:220:GLN:HB2	2.50	0.47
1:C:256:GLU:CD	1:C:283:ARG:HH21	2.19	0.46
1:C:41:ARG:HD2	1:C:77:VAL:CG2	2.37	0.46
1:E:198:GLY:HA2	1:E:222:LEU:HD12	1.96	0.46
1:F:44:ILE:HD12	1:F:54:LEU:HD21	1.96	0.46
1:F:179:LEU:O	1:F:181:PRO:HD3	2.16	0.46
1:C:37:VAL:CG1	1:C:40:ILE:HG13	2.45	0.46
1:B:191:LYS:HD3	3:B:506:HOH:O	2.16	0.46
1:B:304:ASP:O	1:B:306:SER:N	2.39	0.46
1:C:71:ARG:HD3	1:C:103:ASP:OD1	2.16	0.46
1:B:43:ARG:O	1:B:47:VAL:HG23	2.16	0.46
1:F:68:GLU:H	1:F:68:GLU:CD	2.19	0.46
1:C:121:PRO:HA	1:C:154:LEU:HB2	1.98	0.46
1:C:244:GLY:HA2	1:C:271:TRP:CD1	2.51	0.46
1:E:41:ARG:CZ	1:E:73:GLY:CA	2.94	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:PRO:HG3	1:A:166:ALA:HB1	1.97	0.46
1:B:71:ARG:HD3	1:B:103:ASP:HA	1.97	0.46
1:E:43:ARG:HD3	1:E:46:GLU:OE1	2.16	0.46
1:F:337:LYS:O	1:F:339:PRO:HD3	2.16	0.46
1:E:36:HIS:HA	1:E:41:ARG:HH21	1.81	0.46
1:C:154:LEU:HD21	1:C:271:TRP:CZ2	2.51	0.45
1:E:337:LYS:O	1:E:337:LYS:HG3	2.16	0.45
1:F:41:ARG:HG3	1:F:77:VAL:HG23	1.99	0.45
1:B:201:PHE:HB3	1:E:260:ARG:HE	1.82	0.45
1:E:81:THR:O	1:E:114:LYS:NZ	2.43	0.45
1:F:41:ARG:HG3	1:F:77:VAL:CG2	2.46	0.45
1:A:193:THR:HG21	1:A:200:PRO:CD	2.45	0.45
1:A:71:ARG:HD3	1:A:103:ASP:HA	1.99	0.45
1:A:203:GLY:HA2	1:A:220:GLN:HG2	1.99	0.44
1:F:182:ALA:HB2	1:F:195:VAL:HG12	2.00	0.44
1:A:40:ILE:HG23	3:A:413:HOH:O	2.17	0.44
1:D:124:GLU:HG2	1:D:157:THR:OG1	2.18	0.44
1:D:191:LYS:NZ	3:D:407:HOH:O	2.49	0.44
1:A:198:GLY:HA2	1:A:222:LEU:HD12	1.98	0.44
1:A:310:MET:HE3	1:A:310:MET:HB2	1.85	0.44
1:D:305:GLY:HA2	1:D:321:TRP:CD1	2.53	0.44
1:D:227:ILE:O	1:D:231:VAL:HG23	2.19	0.43
1:E:21:GLU:CD	1:E:22:LEU:H	2.21	0.43
1:B:297:ARG:NH2	1:B:312:ASP:OD2	2.43	0.43
1:E:117:VAL:HG22	1:E:150:LEU:HB2	1.99	0.43
1:F:168:ALA:O	3:F:401:HOH:O	2.20	0.43
1:B:164:LEU:HA	1:B:164:LEU:HD12	1.78	0.43
1:B:19:ILE:HD12	1:B:21:GLU:O	2.19	0.43
1:B:6:HIS:HA	1:B:9:LYS:HE3	2.01	0.43
1:B:1:MET:O	1:B:1:MET:HG2	2.19	0.42
1:B:78:ALA:HB3	1:B:111:VAL:HG11	2.00	0.42
1:D:126:PRO:HG3	1:D:166:ALA:HB1	2.01	0.42
1:E:78:ALA:HB3	1:E:111:VAL:HG11	2.00	0.42
1:E:41:ARG:CZ	1:E:73:GLY:HA2	2.49	0.42
1:B:31:HIS:HB3	1:B:54:LEU:HD12	2.01	0.42
1:E:288:ASP:CB	1:E:291:GLU:HG3	2.48	0.42
1:C:22:LEU:HD21	1:C:175:HIS:HD2	1.83	0.42
1:D:59:GLU:HG3	1:D:93:ASN:HA	2.01	0.42
1:C:317:ASP:OD1	1:C:320:THR:HG23	2.20	0.42
1:E:193:THR:CG2	1:E:200:PRO:HD2	2.47	0.42
1:A:40:ILE:O	1:A:44:ILE:HG12	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:209:GLN:O	1:B:213:GLU:N	2.53	0.42
1:B:75:ILE:O	1:B:79:ARG:HG3	2.19	0.42
1:A:322:LYS:O	1:A:326:VAL:HG23	2.20	0.42
1:C:89:TRP:CD1	1:C:117:VAL:HB	2.55	0.42
1:E:153:ALA:O	1:E:177:PHE:HA	2.20	0.42
1:F:328:VAL:O	1:F:332:ARG:HG3	2.19	0.41
1:A:311:ILE:HD12	1:A:316:GLN:HG3	2.02	0.41
1:E:179:LEU:O	1:E:244:GLY:HA3	2.20	0.41
1:D:260:ARG:O	1:D:264:LEU:HG	2.21	0.41
1:B:306:SER:OG	1:B:307:GLY:N	2.54	0.41
1:E:39:LYS:HE3	1:E:40:ILE:CD1	2.50	0.41
1:F:34:PRO:HA	1:F:35:PRO:HD3	1.96	0.41
1:B:71:ARG:HD3	1:B:103:ASP:OD1	2.21	0.41
1:D:19:ILE:HD12	1:D:23:PRO:HD3	2.02	0.41
1:E:35:PRO:O	1:E:41:ARG:NH2	2.52	0.41
1:D:31:HIS:HD1	1:D:272:SER:HG	1.64	0.41
1:D:71:ARG:HD3	1:D:103:ASP:OD1	2.21	0.41
1:F:155:LEU:HD23	1:F:155:LEU:HA	1.95	0.40
1:C:19:ILE:HD13	1:C:19:ILE:HA	1.90	0.40
1:C:269:GLY:HA2	1:C:285:PHE:CG	2.56	0.40
1:D:202:TYR:CD2	1:D:222:LEU:HD23	2.56	0.40
1:A:34:PRO:HA	1:A:35:PRO:HD3	1.93	0.40
1:C:41:ARG:NE	1:C:76:GLU:OE2	2.55	0.40
1:E:164:LEU:HD12	1:E:164:LEU:HA	1.78	0.40
1:E:243:TYR:O	1:E:270:ALA:HA	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	340/347 (98%)	328 (96%)	12 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	340/347 (98%)	326 (96%)	14 (4%)	0	100	100
1	C	337/347 (97%)	324 (96%)	13 (4%)	0	100	100
1	D	341/347 (98%)	333 (98%)	8 (2%)	0	100	100
1	E	342/347 (99%)	333 (97%)	9 (3%)	0	100	100
1	F	336/347 (97%)	327 (97%)	9 (3%)	0	100	100
All	All	2036/2082 (98%)	1971 (97%)	65 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	273/276 (99%)	271 (99%)	2 (1%)	84	93
1	B	274/276 (99%)	270 (98%)	4 (2%)	65	83
1	C	271/276 (98%)	270 (100%)	1 (0%)	91	97
1	D	275/276 (100%)	274 (100%)	1 (0%)	91	97
1	E	275/276 (100%)	272 (99%)	3 (1%)	73	88
1	F	272/276 (99%)	268 (98%)	4 (2%)	65	83
All	All	1640/1656 (99%)	1625 (99%)	15 (1%)	78	91

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	ARG
1	A	194	ARG
1	B	46	GLU
1	B	194	ARG
1	B	308	VAL
1	B	310	MET
1	C	59	GLU
1	D	65	ASP

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Mol	Chain	Res	Type
1	E	39	LYS
1	E	59	GLU
1	E	318	ASP
1	F	41	ARG
1	F	193	THR
1	F	215	ARG
1	F	318	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	TRS	B	401	-	7,7,7	0.44	0	9,9,9	1.06	0
2	TRS	C	401	-	7,7,7	1.36	0	9,9,9	4.93	5 (55%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	TRS	B	401	-	-	0/9/9/9	-
2	TRS	C	401	-	-	7/9/9/9	-

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	TRS	C2-C-N	-9.86	78.54	107.98
2	C	401	TRS	C1-C-N	-6.88	87.44	107.98
2	C	401	TRS	C3-C-N	-6.76	87.79	107.98
2	C	401	TRS	C3-C-C2	3.49	121.65	110.81
2	C	401	TRS	C2-C-C1	3.06	120.31	110.81

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	401	TRS	N-C-C1-O1
2	C	401	TRS	C1-C-C2-O2
2	C	401	TRS	N-C-C2-O2
2	C	401	TRS	C2-C-C3-O3
2	C	401	TRS	N-C-C3-O3
2	C	401	TRS	C3-C-C1-O1
2	C	401	TRS	C2-C-C1-O1

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	344/347 (99%)	-0.13	8 (2%) 60 64	32, 46, 67, 84	0
1	B	344/347 (99%)	0.10	22 (6%) 19 20	33, 48, 76, 95	0
1	C	341/347 (98%)	0.04	11 (3%) 47 52	34, 47, 68, 81	0
1	D	345/347 (99%)	-0.08	5 (1%) 75 78	35, 49, 64, 87	0
1	E	346/347 (99%)	-0.11	9 (2%) 56 59	34, 50, 65, 94	0
1	F	342/347 (98%)	-0.11	8 (2%) 60 64	33, 49, 69, 100	0
All	All	2062/2082 (99%)	-0.05	63 (3%) 49 53	32, 48, 69, 100	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	306	SER	6.0
1	B	304	ASP	5.3
1	B	305	GLY	5.3
1	C	307	GLY	5.2
1	B	303	PRO	5.0
1	A	209	GLN	5.0
1	E	196	GLY	4.5
1	A	214	HIS	4.2
1	E	210	GLU	4.0
1	E	211	GLY	3.7
1	B	15	ALA	3.6
1	F	347	PHE	3.5
1	C	308	VAL	3.5
1	B	300	ASP	3.5
1	D	209	GLN	3.5
1	F	214	HIS	3.4
1	D	213	GLU	3.3
1	B	214	HIS	3.3
1	B	1	MET	3.2

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Mol	Chain	Res	Type	RSRZ
1	E	209	GLN	3.2
1	C	12	ALA	3.1
1	A	304	ASP	2.9
1	A	213	GLU	2.9
1	D	311	ILE	2.8
1	B	13	ALA	2.8
1	D	212	GLN	2.7
1	B	338	ASP	2.6
1	C	15	ALA	2.6
1	B	213	GLU	2.6
1	E	14	GLY	2.6
1	F	82	ASP	2.6
1	B	306	SER	2.5
1	C	13	ALA	2.5
1	E	347	PHE	2.5
1	E	13	ALA	2.5
1	B	308	VAL	2.5
1	C	347	PHE	2.5
1	F	339	PRO	2.5
1	B	12	ALA	2.4
1	B	209	GLN	2.4
1	F	17	GLN	2.4
1	F	1	MET	2.4
1	A	321	TRP	2.4
1	B	307	GLY	2.3
1	B	339	PRO	2.3
1	A	347	PHE	2.3
1	E	212	GLN	2.3
1	A	318	ASP	2.3
1	B	310	MET	2.3
1	C	146	GLN	2.3
1	B	144	GLN	2.2
1	F	208	PRO	2.2
1	C	305	GLY	2.2
1	E	289	VAL	2.2
1	B	343	GLN	2.2
1	D	158	ALA	2.2
1	A	303	PRO	2.2
1	B	302	MET	2.1
1	B	7	PHE	2.1
1	B	18	PRO	2.1
1	F	341	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	75	ILE	2.0
1	C	302	MET	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	TRS	C	401	8/8	0.97	0.14	37,40,42,45	0
2	TRS	B	401	8/8	0.98	0.15	35,36,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.